

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX041719\
 Data File : VX009012.D
 Acq On : 17 Apr 2019 14:14
 Operator : JC/SP
 Sample : K2387-07DL 20X
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KY063MW003-190410DL

Quant Time: Apr 18 02:11:49 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X041519W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 17 06:47:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.66	168	261311	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.86	114	417312	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.11	117	359110	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.08	152	154168	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.06	65	169325	45.98	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.96%	
35) Dibromofluoromethane	5.50	113	124249	46.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.78%	
50) Toluene-d8	8.71	98	519868	48.67	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.34%	
62) 4-Bromofluorobenzene	11.13	95	173285	45.96	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.92%	
Target Compounds						
27) cis-1,2-Dichloroethene	4.60	96	34288	10.104	ug/l	90
44) Trichloroethene	7.21	130	33186	10.638	ug/l	99
64) Tetrachloroethene	9.34	164	113465	41.591	ug/l	97

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX041719\
Data File : VX009012.D
Acq On : 17 Apr 2019 14:14
Operator : JC/SP
Sample : K2387-07DL 20X
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
KY063MW003-190410DL

Quant Time: Apr 18 02:11:49 2019
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X041519W.M
Quant Title : SW846 8260
QLast Update : Wed Apr 17 06:47:58 2019
Response via : Initial Calibration

